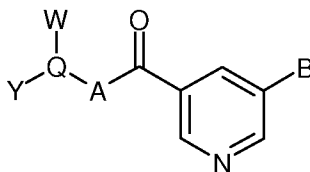


CLAIM AMENDMENTS

Please amend the claims as follows:

1. (currently amended): A compound of formula I



I

or pharmaceutically acceptable ~~prodrugs, salts, hydrates, solvates, crystal forms or~~ diastereomers thereof, wherein:

A is ~~selected from O, S, and~~ NR¹, where R¹ is H, or C₁₋₄ alkyl;

B is ~~aryl, or hetaryl, phenyl~~ optionally substituted with 0-3 ~~with 0-4~~ substituents independently selected from halogen, C₁₋₄ alkyl, CF₃, CN, aryl, ~~hetaryl~~, OH, OCF₃, OC₁₋₄ alkyl, OC₂₋₅ alkylNR²R³, Oaryl, ~~Ohetaryl~~, CO₂R², CONR²R³, NR²R³, ~~C₁₋₄ alkyl~~NR²R³, NR⁴C₁₋₄ alkylNR²R³, NR²COR³, OC(O)NR²R³, NR⁴CONR²R³, and NR²SO₂R³;

wherein R², R³ are each independently H, C₁₋₄ alkyl, ~~C₁₋₄ alkyl~~ heterocyclyl, aryl, ~~hetaryl~~, C₁₋₄ alkyl aryl, ~~C₁₋₄ alkyl hetaryl~~, or may be joined to form an optionally substituted 3-8 membered ring optionally containing one of O, S, or NR⁵;

wherein R⁴ is H or C₁₋₄ alkyl; and

wherein R⁵ is H or C₁₋₄ alkyl;

Q is a bond when W is absent, and is C₁₋₄ alkyl when W is present;

W is selected from H, C₁₋₄ alkyl, C₂₋₆ alkenyl; where C₁₋₄ alkyl or C₂₋₆ alkenyl may be optionally substituted with C₁₋₄ alkyl, OH, OC₁₋₄ alkyl, NR⁶C(O)R⁷, CONR⁶R⁷, OR⁶, or NR⁶R⁷;

wherein R⁶, and R⁷ are each independently H, C₁₋₄ alkyl, C₁₋₄ alkyl cycloalkyl, ~~C₁₋₄ alkyl~~ heterocyclyl, aryl, ~~hetaryl~~, or may be joined to form an optionally substituted 3-8 membered ring ~~optionally containing one of O, S, or NR⁸~~, and

~~wherein R⁸ is H, or C₁₋₄ alkyl;~~

Y is H ~~or phenyl~~, aryl ~~or hetaryl~~ optionally substituted with 0-3 substituents independently selected from halogen, C₁₋₄ alkyl, CF₃, aryl, ~~hetaryl~~, OH, OCF₃, CN, C₂₋₄ alkynyl, OC₁₋₄ alkyl,

OC₂₋₅ alkylNR⁹R¹⁰, Oaryl, ~~Ohetaryl~~, CO₂R⁹, CONR⁹R¹⁰, NR⁹R¹⁰, C₁₋₄ alkylNR⁹R¹⁰,
NR¹¹C₁₋₄ alkylNR⁹R¹⁰, NR⁹COR¹⁰, NR¹¹CONR⁹R¹⁰, and NR⁹SO₂R¹⁰;

wherein R⁹ and R¹⁰ is each independently H, C₁₋₄ alkyl, ~~C₁₋₄ alkyl heterocyclyl~~, aryl, ~~hetaryl~~,
C₁₋₄ alkyl aryl, ~~C₁₋₄ alkyl hetaryl~~, or may be joined to form an optionally substituted
3-8 membered ring ~~optionally containing one O, S, or NR¹²~~;

wherein R¹¹ is H or C₁₋₄ alkyl; ~~and R¹² is H or C₁₋₄ alkyl~~

with the proviso that when Y is phenyl substituted at the ortho position with CO₂R⁹, CN or
NH₂, W is absent and Q is a bond, and B has one or zero substituents, then R¹ is C₁₋₄ alkyl.

2. (currently amended): A compound according to claim 1 or pharmaceutically acceptable ~~prodrugs, salts, hydrates, solvates, crystal forms~~ or diastereomers thereof, wherein:

A is NR¹ and R¹ is H or C₁₋₄ alkyl;

W is selected from H, C₁₋₄ alkyl, and C₂₋₆ alkenyl; ~~where~~ wherein C₁₋₄ alkyl or C₂₋₆ alkenyl may be optionally substituted with C₁₋₄ alkyl, OH, OC₁₋₄ alkyl, or NR⁶R⁷;

wherein R⁶, and R⁷ are each independently H, C₁₋₄ alkyl, C₁₋₄ alkyl cycloalkyl, ~~C₁₋₄ alkyl heterocyclyl~~, aryl, ~~hetaryl~~, or may be joined to form an optionally substituted 3-8 membered ring ~~optionally containing one of O, S or NR⁸~~;

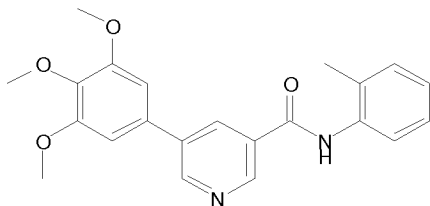
~~wherein R⁸ is H or C₁₋₄ alkyl;~~

Y is H or phenyl, aryl or ~~hetaryl~~ optionally substituted with 0-3 substituents independently selected from halogen, C₁₋₄ alkyl, CF₃, aryl, ~~hetaryl~~, OH, OCF₃, OC₁₋₄ alkyl, OC₂₋₅ alkylNR⁹R¹⁰, Oaryl, ~~Ohetaryl~~, CO₂R⁹, CONR⁹R¹⁰, NR⁹R¹⁰, C₁₋₄ alkylNR⁹R¹⁰, NR¹¹C₁₋₄ alkylNR⁹R¹⁰, NR⁹COR¹⁰, NR¹¹CONR⁹R¹⁰, and NR⁹SO₂R¹⁰;

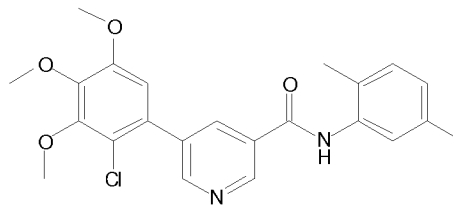
wherein R⁹, and R¹⁰ are each independently H, C₁₋₄ alkyl, ~~C₁₋₄ alkyl heterocyclyl~~, aryl, ~~hetaryl~~, C₁₋₄ alkyl aryl, ~~C₁₋₄ alkyl hetaryl~~, or may be joined to form an optionally substituted 3-8 membered ring ~~optionally containing an atom selected from one of O, S, NR¹²~~; and ~~R¹¹ is selected from H, or NR¹²~~;

wherein R¹¹ is H or C₁₋₄ alkyl; ~~and R¹² is selected from H, R¹² is H or C₁₋₄ alkyl.~~

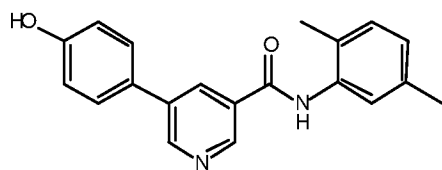
3. (currently amended): A compound according to claim 1 wherein the compound is selected from the group consisting of:



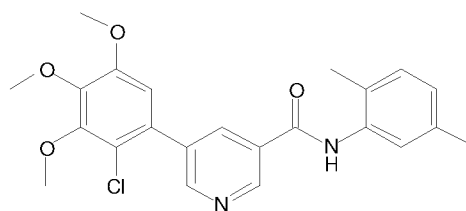
C22H22N2O4



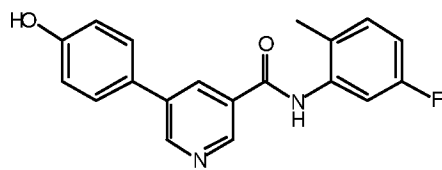
C23H23ClN2O4



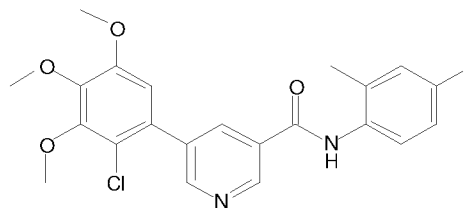
C20H18N2O2



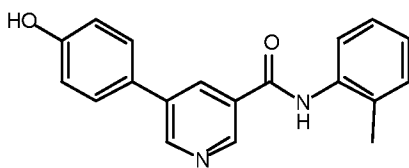
C22H20ClFN2O4



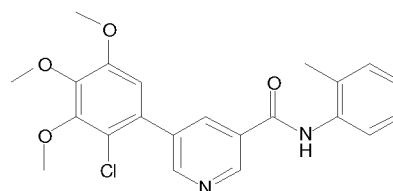
C19H15FN2O2



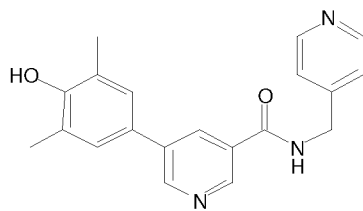
C22H20ClFN2O4



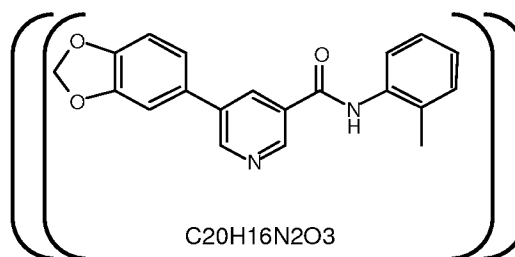
C19H16N2O2



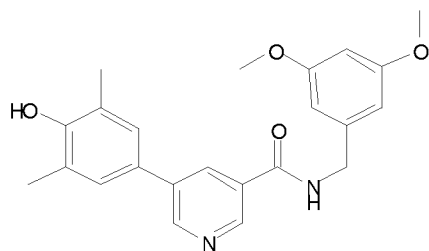
C22H21ClN2O4



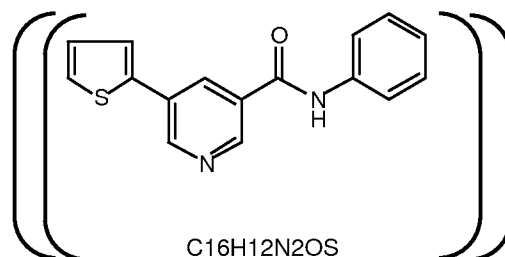
C20H19N3O2



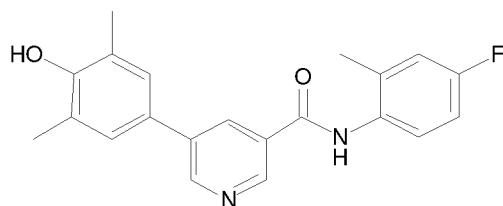
C20H16N2O3



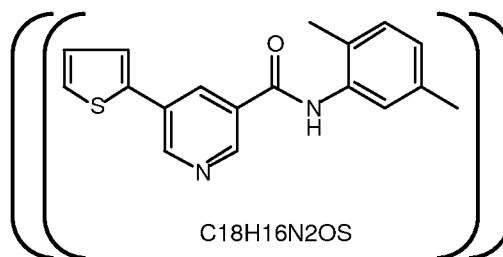
C23H24N2O4



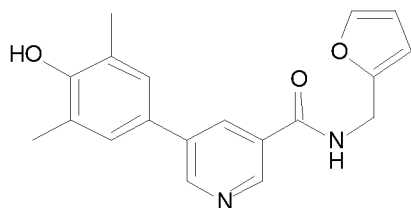
C16H12N2OS



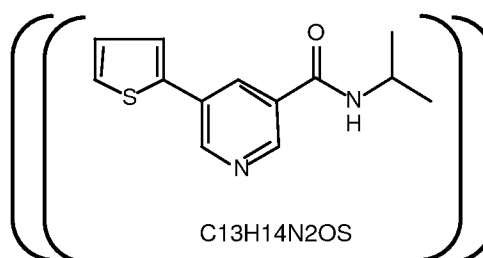
C21H19FN2O2



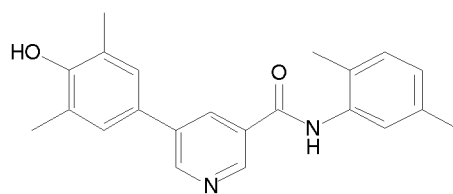
C18H16N2OS



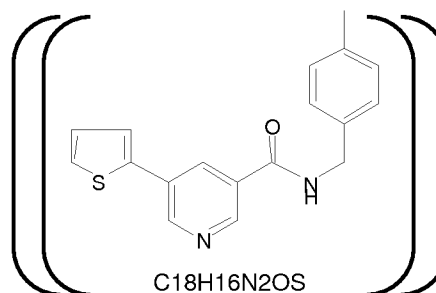
C19H18N2O3



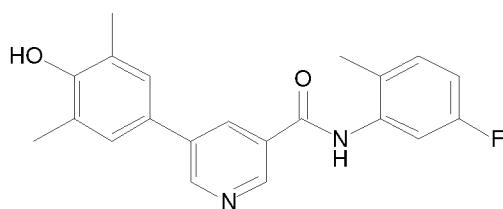
C13H14N2OS



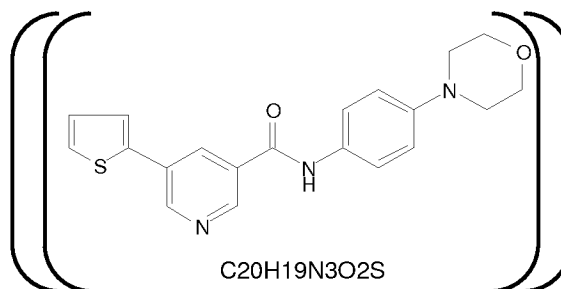
C22H22N2O2



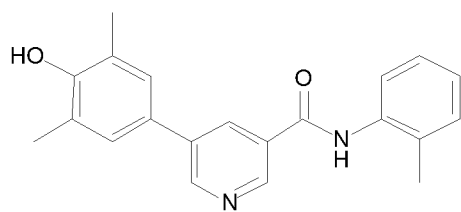
C18H16N2OS



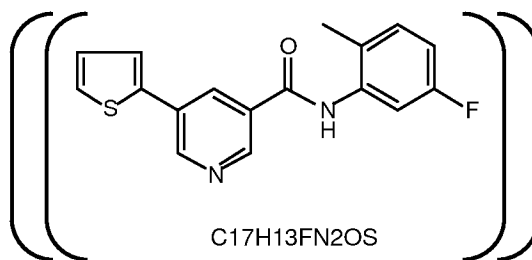
C21H19FN2O2



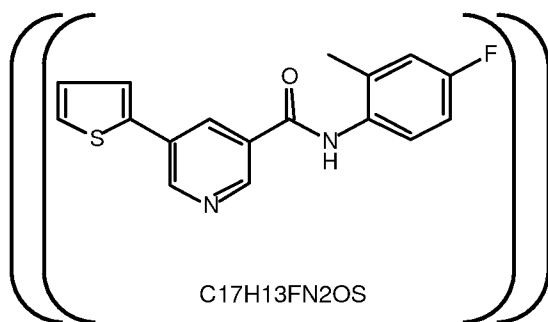
C20H19N3O2S



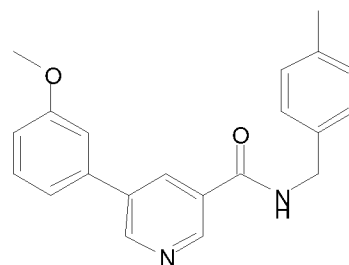
C21H20N2O2



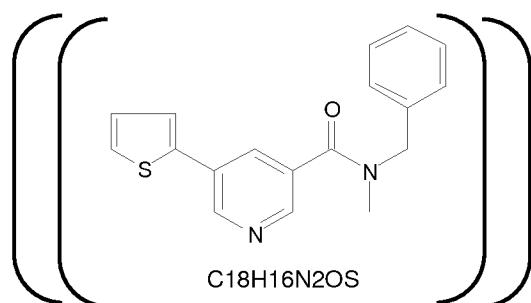
C17H13FN2OS



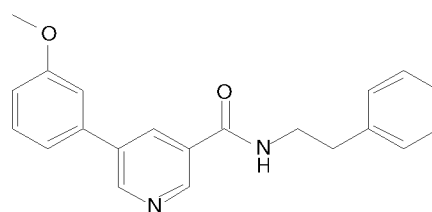
C17H13FN2OS



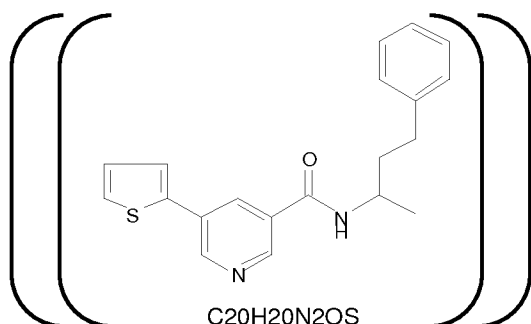
C21H20N2O2



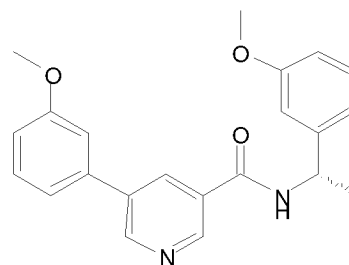
C18H16N2OS



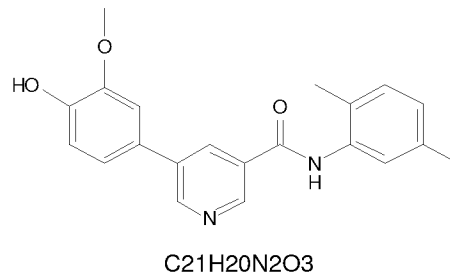
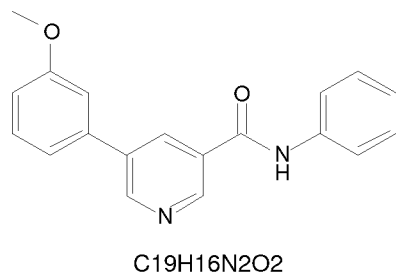
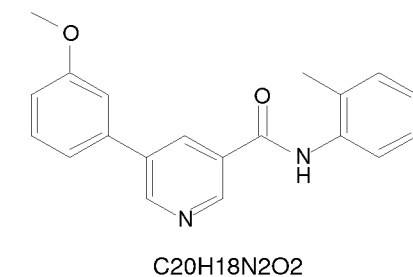
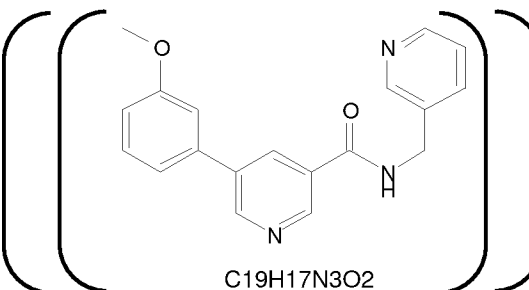
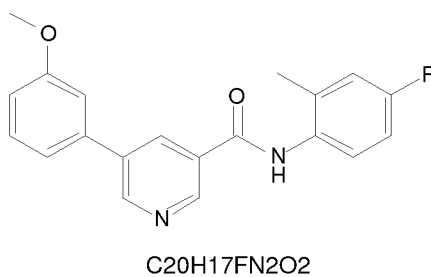
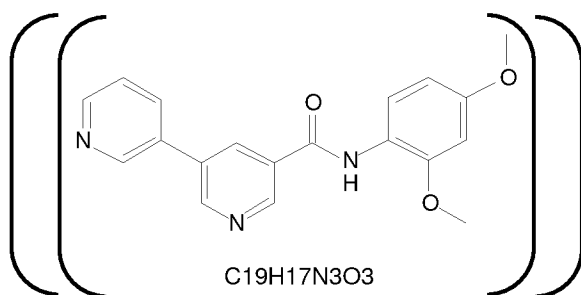
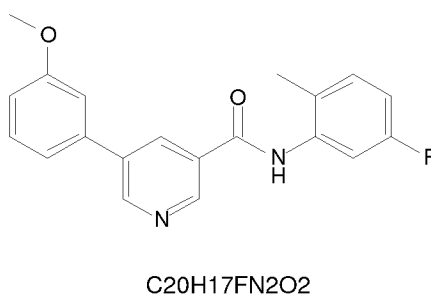
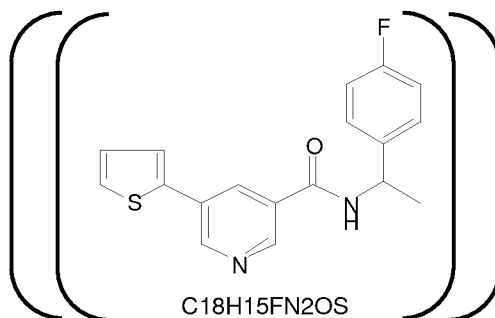
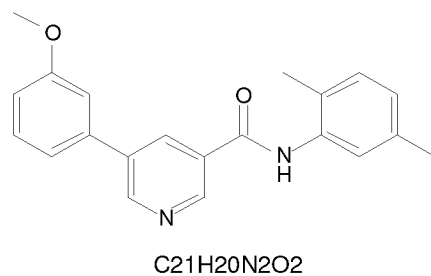
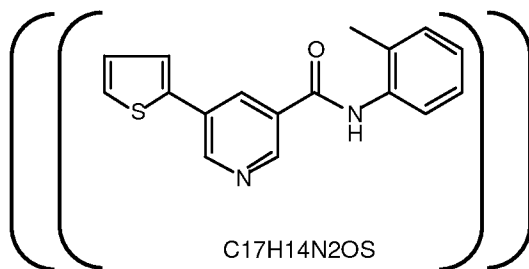
C21H20N2O2

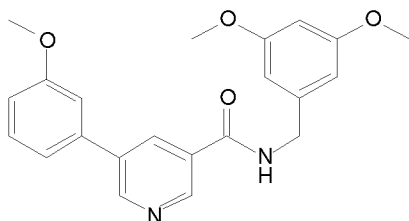


C20H20N2OS

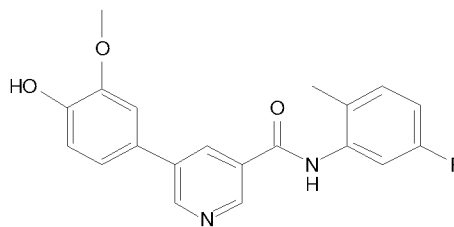


C22H22N2O3

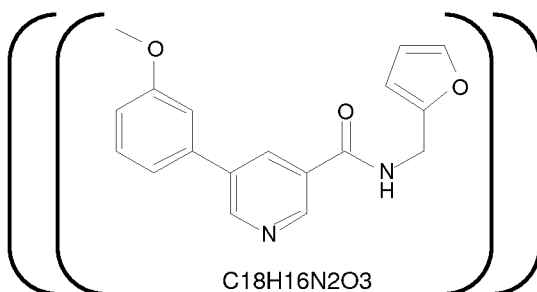




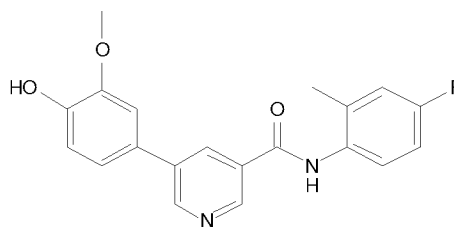
C22H22N2O4



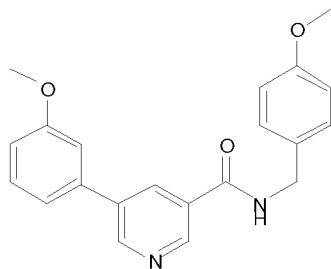
C20H17FN2O3



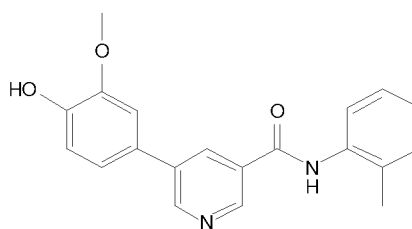
C18H16N2O3



C20H17FN2O3



C21H20N2O3



C20H18N2O3

or pharmaceutically acceptable prodrugs, salts, hydrates, solvates, crystal forms or diastereomers thereof.

4. (currently amended): A pharmaceutical composition comprising a carrier and at least one compound of claim 1.

5-6. (canceled)

7. (new): A pharmaceutical composition comprising a carrier and at least one compound of claim 2.

8. (new): A pharmaceutical composition comprising a carrier and at least one compound of claim 3.